

a compound having the formula:



*32* or a pharmaceutically acceptable salt or hydrate thereof, wherein:

$n$  is 0, 1, 2, 3 or 4;

$X$  is absent, ( $\text{C}_1\text{-C}_3$ ) alkyl, ( $\text{C}_1\text{-C}_3$ ) alkenyl, or ( $\text{C}_1\text{-C}_3$ ) alkynyl;

$Y$  is C, N, P, Si or Ge;

$\text{R}_1$  is absent, -halo, -R, -OR, -SR, -NR<sub>2</sub>, -ONR<sub>2</sub>, -NO<sub>2</sub>, -CN, -C(O)R, -C(S)R, -C(O)OR, -C(S)OR, -C(O)SR, -C(S)SR, -C(O)NR<sub>2</sub>, -C(S)NR<sub>2</sub>, -C(O)NR(OR), -C(S)NR(OR), -C(O)NR(SR), C(S)NR(SR), -CH(CN)<sub>2</sub>, -CH[C(O)R]<sub>2</sub>, -CH[C(S)R]<sub>2</sub>, -CH[C(O)OR]<sub>2</sub>, -CH[C(S)OR]<sub>2</sub>, -CH[C(O)SR]<sub>2</sub>, -CH[C(S)SR]<sub>2</sub> or aryl;

$\text{Ar}_1$  is aryl, substituted aryl, heteroaryl other than imidazole, nitroimidazole and triazole, heteroarylium other than imidazolium, nitroimidazolium and triazolium, ( $\text{C}_5\text{-C}_8$ ) cycloalkyl or ( $\text{C}_5\text{-C}_8$ ) heterocycloalkyl;

$\text{Ar}_2$  is aryl or substituted aryl;

$\text{Ar}_3$  is aryl, substituted aryl, biaryl or heteroaryl other than imidazole, nitroimidazole and triazole; each R is independently selected from the group consisting of -H, ( $\text{C}_1\text{-C}_6$ ) alkyl, substituted ( $\text{C}_1\text{-C}_6$ ) alkyl, ( $\text{C}_1\text{-C}_6$ ) alkenyl, substituted ( $\text{C}_1\text{-C}_6$ ) alkenyl ( $\text{C}_1\text{-C}_6$ ) alkynyl, substituted ( $\text{C}_1\text{-C}_6$ ) alkynyl, and ( $\text{C}_1\text{-C}_6$ ) alkoxy;

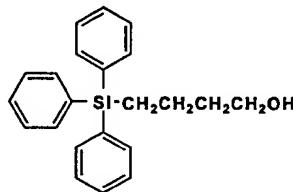
the aryl substituents are each independently selected from the group consisting of -halo, trihalomethyl, -R, -R', -OR', -SR', NR'<sub>2</sub>, -NO<sub>2</sub>, -CN, -C(O)R', -C(S)R', -C(O)OR', -C(S)OR', -C(O)SR' and -C(S)SR';

the alkyl, alkenyl and alkynyl substituents are each independently selected from the group consisting of -halo, -R', -OR', -SR', NR'<sub>2</sub>, -NO<sub>2</sub>, -CN, -C(O)R', -C(S)R', -C(O)OR', -C(S)OR', -C(O)SR', -C(S)SR', aryl,  $\gamma$ -butyrolactonyl, pyrrolidinyl and succinic anhydridyl; [and]

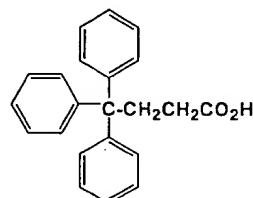
each R' is independently selected from the group consisting of -H, ( $\text{C}_1\text{-C}_6$ ) alkyl, ( $\text{C}_1\text{-C}_6$ ) alkenyl and ( $\text{C}_1\text{-C}_6$ ) alkynyl.

TABLE A  
Exemplary Compounds

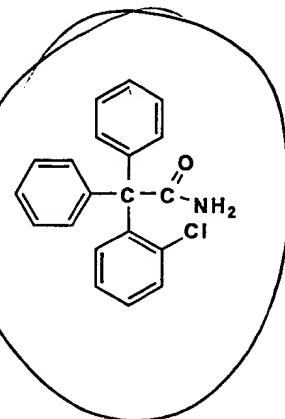
Chemical Exemplar



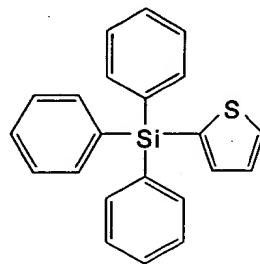
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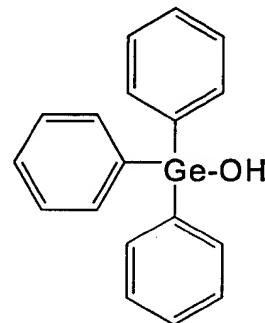
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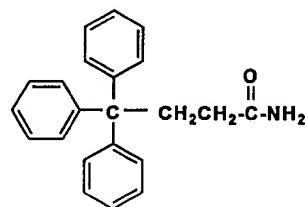
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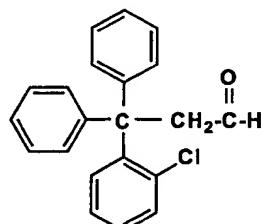
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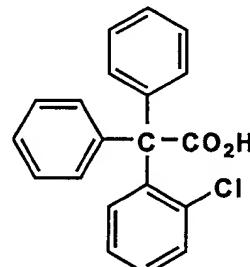
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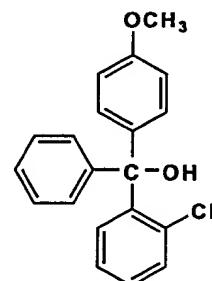
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